WHAT IS CLAIMED IS:

1. An indene derivative of formula (I) or a pharmaceutically acceptable salt thereof:

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(I)

wherein,

 R_{1a} is OH or H;

 R_{1b} is C_{1-6} alkyl, C_{3-6} cycloalkyl, benzyl or phenyl, the phenyl being optionally substituted with one or more substituents selected from the group consisting of halogen, CN, NH_2 , NO_2 and OR^a , when R_{1a} is OH; when R_{1a} is H,

$$R_{1b}$$
 is OR^a , NR^bR^c , $NHCOR^a$ or

R₂ is CN, CO₂R^a or CONR^eR^f;

R₃ is phenyl optionally substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂, NO₂, OR^a and C₁₋₆ alkyl; and

 $R^4,\,R^5,\,R^6$ and R^7 are each independently H, $O(CH_2)_mR^g$ or $CH_2R^h;$ in which

 R^a is H, C_{1-6} alkyl or C_{3-6} cycloalkyl, the C_{1-6} alkyl and C_{3-6} cycloalkyl being optionally substituted with one or more halogens;

 $R^b,\,R^c,\,R^e$ and R^f are each independently H, $C_{1\text{-}6}$ alkyl, $C_{3\text{-}6}$ cycloalkyl or benzyl;

R^d is O, S or NR^a;

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 R^g is H, , or phenyl, the phenyl being optionally substituted with one or more substituents selected from the group consisting of halogen, CN, NH₂ and NO₂;

$$R_h$$
 is $-\xi-N$ R^d ; and

m is an integer in the range of 1 to 3.

2. The compound of claim 1, wherein R_{1b} is C_{1-6} alkyl, C_{3-6} cycloalkyl, benzyl or phenyl, the phenyl being optionally substituted with one or more methoxy groups,

when R_{1a} is OH; when R_{1a} is H, R_{1b} is OR^a, NR^bR^c, NHCOR^a or R_{1a} ; R₃ is phenyl being optionally substituted with one or more halogens or R_{1a} alkyls; and R_{1a} and R_{1a} is H, in which R_{1a} is H or R_{1a} alkyl; R_{1a} is O or S; R_{1a} is H, phenyl,

$$-\frac{1}{2} \frac{1}{U_N} \qquad -\frac{1}{2} \frac{1}{U_N} \qquad \text{or} \qquad R^d$$

- 3. The compound of claim 1, wherein R_3 is phenyl, R_5 is H, and R_6 is $O(CH_2)_m R^g$ or CH_2R^h .
 - 4. The compound of claim 1, which is selected from the group consisting of:

 1-hydroxy-6-methoxy-1,3-diphenyl-1H-indene-2-carboxylic acid ethyl ester,
 - 1-hydroxy-6-methoxy-1-(3-methoxy-phenyl)-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-1-isopropyl-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-6-methoxy-1-methyl-3-phenyl-1H-indene-2-carboxylic acid

ethyl ester,

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1-benzyl-1-hydroxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-cyclohexyl-1-hydroxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-1,3-diphenyl-6-(3-phenyl-propoxy)-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-6-(2-morpholine-4-yl-ethoxy)-1,3-diphenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-6-morpholine-4-yl-methyl-1,3-diphenyl-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-1,3-diphenyl-6-(2-pyridine-2-yl-ethoxy)-1H-indene-2-carboxylic acid ethyl ester,

1-hydroxy-1,3-diphenyl-6-(3-phenyl-propoxy)-1H-indene-2-Carbonitrile,

1-hydroxy-1,3-diphenyl-6-(3-phenyl-propoxy)-1H-indene-2-carboxylic acid methyl ester,

1-hydroxy-6-methoxy-1,3-diphenyl-1H-indene-2-carboxylic acid,

1-hydroxy-6-methoxy-1-methyl-3-phenyl-1H-indene-2-carboxylic acid, 1-benzyl-1-hydroxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid,

1-hydroxy-1,3-diphenyl-6-(3-phenyl-propoxy)-1H-indene-carboxylic acid,

1-cyclohexyl-1-hydroxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid,

1,6-dimethoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-ethoxy-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-amino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-amino-3-phenyl-6-(3-phenyl-propoxy)-1H-indene-2-carboxylic acid ethyl ester,

1-amino-6-(2-morpholin-4-yl-ethoxy)-3-phenyl-1H-indene-2-carboxylic acid cyclohexyl amide,

1-amino-3-phenyl-6-(3-phenyl-propoxy)-1H-indene-2-carbonitrile,

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1-acetylamino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

6-methoxy-3-phenyl-1-propionylamino-1H-indene-2-carboxylic acid ethyl ester,

1-acetylamino-3-phenyl-6-(3-phenyl-propoxy)-1H-indene-2-carboxylic acid ethyl ester,

1-acetylamino-6-(2-morpholin-4-yl-ethoxy)-3-phenyl-1H-indene-2-carboxylic acid cyclohexyl amide,

1-diethylamino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-ethylamino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester, 6-methoxy-1-morpholin-4-yl-3-phenyl-1H-indene-2-carboxylic acid ethyl ester,

1-benzyl amino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester, and

1-cyclohexyl amino-6-methoxy-3-phenyl-1H-indene-2-carboxylic acid ethyl ester.

5. A process for preparing a compound of formula (I-a) which comprises reacting a compound of formula (II) with a Grignard reagent:

(II)

wherein R_{1a} is OH; R_{1b} is alkyl, phenyl or benzyl; and R₂, R₃, R₄, R₅, R₆ and R₇ have

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the same meaning as defined in claim 1.

6. A process for preparing a compound of formula (I-e) which comprises reacting a compound of formula (II) with hydroxyl amine to obtain a compound of formula (III), and hydrogenation of the compound of formula (III) followed by reacting with acetyl chloride or an anhydrous acetic acid:

wherein R_{1a} is H; R_{1b} is NH_2 or $NHCOR^a$; and R_2 , R_3 , R_4 , R_5 , R_6 and R_7 have the same meaning as defined in claim 1.

7. A process for preparing a compound of formula (I-d) which comprises halogenation of a compound of formula (VIII) to obtain a compound of formula (IV), and reacting the compound of formula (IV) with an amine or alcohol compound:

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$$R_6$$
 R_7
 R_7
 R_8
 R_7
 R_8
 R_8
 R_9
 R_9

wherein R_{1a} is H; R_{1b} is OR^a , NR^bR^c or R_6 and R_7 have the same meaning as defined in claim 1.

- 8. A pharmaceutical composition for modulating the activities of peroxisome proliferator activated receptors (PPARs) comprising a therapeutically effective amount of the compound or a salt defined in claim 1 as an active ingredient together with a pharmaceutically acceptable carrier.
- 9. The composition of claim 8, which is used for the treatment and prevention of diabetes, obesity, arteriosclerosis, hyperlipidemia, hyperinsulinism, hypertension, osteoporosis, liver cirrhosis, asthma and cancer.

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